



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2026 – 08:12 AM EDT

PDB ID : 7T97 / pdb_00007t97
Title : Crystal structure of engineered CYS-CYS fab dimer CH1-207 (HC4)
Authors : Harris, S.F.; Boenig, G.D.L.
Deposited on : 2021-12-18
Resolution : 2.14 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

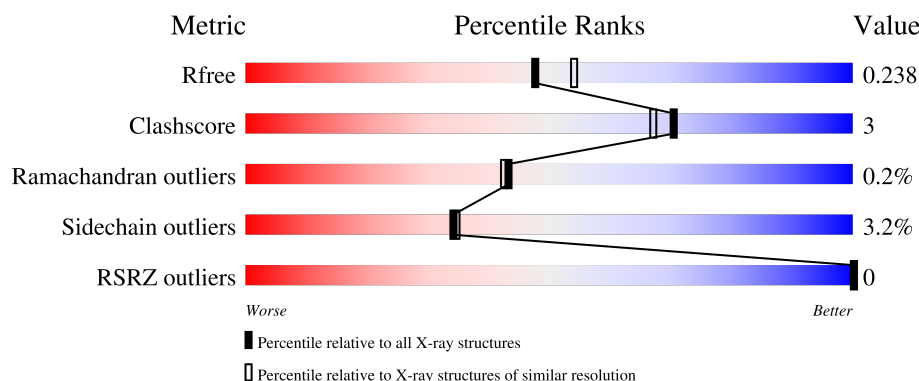
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.14 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3689 (2.16-2.12)
Clashscore	190562	3812 (2.16-2.12)
Ramachandran outliers	187476	3773 (2.16-2.12)
Sidechain outliers	187428	3772 (2.16-2.12)
RSRZ outliers	180081	3691 (2.16-2.12)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	228	
1	C	228	
1	E	228	
1	H	228	
2	B	214	

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Mol	Chain	Length	Quality of chain
2	D	214	88% 11% •
2	F	214	89% 10% •
2	L	214	91% 7% ••

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 13931 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called FAB Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	220	Total	C	N	O	S	0	3	0
			1657	1048	279	320	10			
1	C	221	Total	C	N	O	S	0	3	0
			1666	1053	280	323	10			
1	E	220	Total	C	N	O	S	0	3	0
			1657	1048	279	320	10			
1	H	220	Total	C	N	O	S	0	3	0
			1657	1048	279	320	10			

- Molecule 2 is a protein called FAB Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	212	Total	C	N	O	S	0	1	0
			1637	1026	275	330	6			
2	D	212	Total	C	N	O	S	0	1	0
			1637	1026	275	331	5			
2	F	212	Total	C	N	O	S	0	2	0
			1640	1028	275	331	6			
2	L	212	Total	C	N	O	S	0	2	0
			1640	1028	275	331	6			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	92	Total	O	0	0
			92	92		
3	B	86	Total	O	0	0
			86	86		
3	C	99	Total	O	0	0
			99	99		
3	D	90	Total	O	0	0
			90	90		

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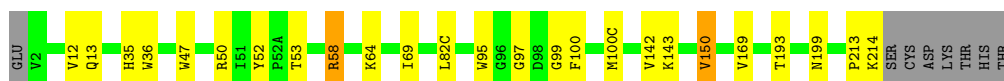
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	84	Total 84	O 84	0	0
3	F	99	Total 99	O 99	0	0
3	H	77	Total 77	O 77	0	0
3	L	113	Total 113	O 113	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

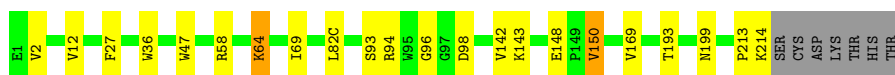
- Molecule 1: FAB Heavy chain

Chain A: 



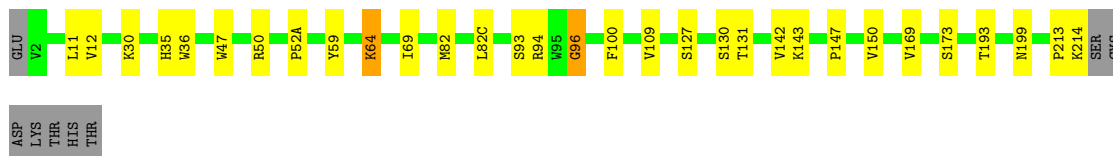
- Molecule 1: FAB Heavy chain

Chain C: 



- Molecule 1: FAB Heavy chain

Chain E: 



- Molecule 1: FAB Heavy chain

Chain H: 



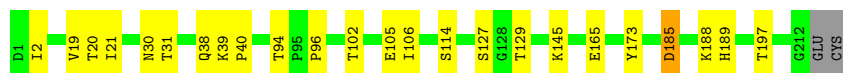
- Molecule 2: FAB Light Chain

Chain B: 



- Molecule 2: FAB Light Chain

Chain D:  88% 11% •



• Molecule 2: FAB Light Chain

Chain F:  89% 10% •



• Molecule 2: FAB Light Chain

Chain L:  91% 7% ••



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	151.38Å 141.07Å 118.61Å 90.00° 90.32° 90.00°	Depositor
Resolution (Å)	75.69 – 2.14 75.69 – 2.14	Depositor EDS
% Data completeness (in resolution range)	66.6 (75.69-2.14) 66.4 (75.69-2.14)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 2.14Å)	Xtriage
Refinement program	BUSTER 2.11.7 (20-MAY-2020)	Depositor
R, R_{free}	0.194 , 0.227 0.207 , 0.238	Depositor DCC
R_{free} test set	4656 reflections (3.42%)	wwPDB-VP
Wilson B-factor (Å ²)	32.6	Xtriage
Anisotropy	0.067	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 22.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.104 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13931	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 44.47 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.5194e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.71	0/1705	0.95	1/2324 (0.0%)
1	C	0.72	0/1714	0.95	1/2336 (0.0%)
1	E	0.73	0/1705	0.95	1/2324 (0.0%)
1	H	0.72	0/1705	0.97	0/2324
2	B	0.69	0/1677	0.99	2/2280 (0.1%)
2	D	0.70	0/1677	0.98	2/2280 (0.1%)
2	F	0.67	0/1683	0.96	1/2288 (0.0%)
2	L	0.71	0/1683	1.01	3/2288 (0.1%)
All	All	0.71	0/13549	0.97	11/18444 (0.1%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	1	ASP	CA-C-N	6.16	129.31	122.22
2	L	1	ASP	C-N-CA	6.16	129.31	122.22
2	D	185	ASP	CA-CB-CG	5.61	118.21	112.60
2	L	114	SER	N-CA-C	-5.42	100.47	109.46
1	A	13	GLN	CB-CG-CD	-5.41	103.40	112.60
2	B	114	SER	N-CA-C	-5.29	100.68	109.46
2	B	30	ASN	CA-CB-CG	5.21	117.81	112.60
2	D	114	SER	N-CA-C	-5.16	100.89	109.46
1	C	143	LYS	N-CA-C	5.14	117.95	109.72
2	F	114	SER	N-CA-C	-5.08	101.03	109.46
1	E	143	LYS	N-CA-C	5.05	117.80	109.72

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1657	0	1618	12	0
1	C	1666	0	1627	9	0
1	E	1657	0	1618	14	0
1	H	1657	0	1618	15	0
2	B	1637	0	1593	7	0
2	D	1637	0	1594	14	0
2	F	1640	0	1598	15	0
2	L	1640	0	1598	10	0
3	A	92	0	0	0	0
3	B	86	0	0	1	0
3	C	99	0	0	0	0
3	D	90	0	0	3	0
3	E	84	0	0	0	0
3	F	99	0	0	9	0
3	H	77	0	0	5	0
3	L	113	0	0	4	0
All	All	13931	0	12864	88	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (88) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:102:THR:HG23	3:L:315:HOH:O	1.42	1.16
1:H:152:VAL:HG13	3:H:324:HOH:O	1.51	1.11
2:F:39:LYS:NZ	3:F:301:HOH:O	1.87	1.04
1:A:58:ARG:HD3	2:B:94:THR:HG21	1.54	0.89
2:B:22:THR:HG22	2:B:72:THR:HG22	1.60	0.80
1:C:36:TRP:HD1	1:C:69:ILE:HD12	1.47	0.78
1:E:96:GLY:HA3	1:E:100:PHE:HB3	1.69	0.75
1:H:36:TRP:HD1	1:H:69:ILE:HD12	1.50	0.74
2:F:211:ARG:HD3	3:F:302:HOH:O	1.87	0.74
1:A:36:TRP:HD1	1:A:69:ILE:HD12	1.54	0.72
1:A:52:TYR:CE2	1:A:53:THR:HB	2.30	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:82:MET:HG3	3:H:302:HOH:O	1.96	0.66
2:D:189:HIS:HD2	3:D:348:HOH:O	1.80	0.64
1:C:36:TRP:CD1	1:C:69:ILE:HD12	2.32	0.64
2:F:186:TYR:O	3:F:302:HOH:O	2.15	0.64
2:L:100:GLN:HG3	3:L:411:HOH:O	1.99	0.63
1:H:36:TRP:HZ3	3:H:301:HOH:O	1.82	0.62
2:D:39:LYS:NZ	3:D:301:HOH:O	2.29	0.61
2:L:2:ILE:H	2:L:2:ILE:HD13	1.65	0.60
2:D:40:PRO:HG2	2:D:165:GLU:HG3	1.84	0.58
2:F:189:HIS:O	3:F:302:HOH:O	2.17	0.58
1:H:36:TRP:CD1	1:H:69:ILE:HD12	2.37	0.58
1:H:213:PRO:O	1:H:214:LYS:HB2	2.03	0.58
2:F:37:GLN:HB2	2:F:47:LEU:HD11	1.86	0.58
1:H:142:VAL:HG11	1:H:150:VAL:HG11	1.85	0.57
1:C:213:PRO:O	1:C:214:LYS:HB2	2.02	0.56
2:F:29:VAL:HB	2:F:92:TYR:HB2	1.87	0.56
1:E:213:PRO:O	1:E:214:LYS:HB2	2.05	0.56
1:A:213:PRO:O	1:A:214:LYS:HB2	2.06	0.56
1:E:59:TYR:HB2	1:E:64:LYS:HE3	1.88	0.55
1:H:6:GLU:HG2	3:H:301:HOH:O	2.07	0.54
1:C:142:VAL:HG11	1:C:150:VAL:HG11	1.89	0.54
1:A:36:TRP:CD1	1:A:69:ILE:HD12	2.41	0.53
1:C:47:TRP:CD2	2:D:96:PRO:HD2	2.44	0.53
1:H:107:THR:HG22	3:H:301:HOH:O	2.08	0.53
1:E:50:ARG:NH1	2:F:94:THR:OG1	2.42	0.53
2:D:185:ASP:HA	2:D:188:LYS:HE2	1.90	0.52
2:L:2:ILE:HG13	2:L:90:GLN:NE2	2.25	0.52
2:F:211:ARG:CD	3:F:302:HOH:O	2.52	0.52
1:E:142:VAL:HG11	1:E:150:VAL:HG11	1.91	0.52
1:A:142:VAL:HG11	1:A:150:VAL:HG11	1.91	0.52
2:B:127:SER:HB2	1:H:173:SER:HA	1.92	0.51
1:E:36:TRP:HD1	1:E:69:ILE:HD12	1.76	0.51
1:E:82:MET:HE1	1:E:109:VAL:HG21	1.92	0.51
2:F:103:LYS:HG3	3:F:307:HOH:O	2.09	0.51
2:B:37:GLN:HB2	2:B:47:LEU:HD11	1.92	0.50
1:E:12:VAL:HG11	1:E:82(C):LEU:HD13	1.94	0.49
1:A:12:VAL:HG11	1:A:82(C):LEU:HD13	1.93	0.49
1:H:11:LEU:HB2	1:H:147:PRO:HG3	1.95	0.49
1:C:58:ARG:HB3	2:D:94:THR:HG21	1.95	0.48
2:L:21:ILE:HG12	2:L:102:THR:HG21	1.95	0.48
2:F:21:ILE:HG12	2:F:102:THR:HG21	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:12:VAL:HG11	1:H:82(C):LEU:HD13	1.95	0.48
1:C:12:VAL:HG11	1:C:82(C):LEU:HD13	1.95	0.48
2:D:127:SER:HA	1:E:173:SER:HA	1.96	0.48
1:A:50:ARG:HH12	2:B:94:THR:HG1	1.60	0.48
2:D:105:GLU:HG3	2:D:173:TYR:OH	2.14	0.47
2:D:105:GLU:HG2	2:D:106:ILE:N	2.30	0.47
2:L:8:PRO:HG2	3:L:315:HOH:O	2.15	0.47
2:D:38:GLN:NE2	3:D:304:HOH:O	2.49	0.47
2:D:21:ILE:HG12	2:D:102:THR:HG21	1.96	0.46
2:L:35:TRP:HB2	2:L:48:ILE:HB	1.97	0.46
2:D:30:ASN:CG	2:D:31:THR:H	2.24	0.45
2:B:21:ILE:HG12	2:B:102:THR:HG21	1.97	0.45
2:L:145:LYS:HB3	2:L:197:THR:HB	1.98	0.45
2:B:145:LYS:HB3	2:B:197:THR:HB	1.99	0.45
2:D:145:LYS:HB3	2:D:197:THR:HB	1.99	0.44
2:L:124:GLN:NE2	3:L:303:HOH:O	2.49	0.44
2:F:145:LYS:HB3	2:F:197:THR:HB	1.99	0.44
1:E:127:SER:O	1:E:131:THR:HG23	2.18	0.43
1:C:2:VAL:HG13	1:C:27:PHE:CD1	2.54	0.43
1:C:64:LYS:HB2	1:C:64:LYS:HE2	1.85	0.43
1:A:143:LYS:HE2	3:B:303:HOH:O	2.19	0.42
2:F:211:ARG:NE	3:F:302:HOH:O	2.51	0.42
1:H:61:ASP:HA	1:H:64:LYS:HG2	2.01	0.42
1:H:51:ILE:O	1:H:52(A):PRO:HD3	2.20	0.42
1:H:130:SER:O	1:H:136:ALA:HA	2.20	0.42
2:F:100:GLN:HG3	3:F:399:HOH:O	2.20	0.42
1:E:30:LYS:HA	1:E:52(A):PRO:HB2	2.01	0.41
2:F:30:ASN:ND2	2:F:31:THR:H	2.18	0.41
2:L:90:GLN:HE21	2:L:97:THR:H	1.68	0.41
1:A:35:HIS:ND1	1:A:47:TRP:NE1	2.65	0.41
2:F:192:TYR:HE1	3:F:302:HOH:O	2.03	0.41
1:E:35:HIS:ND1	1:E:47:TRP:NE1	2.64	0.40
2:D:127:SER:HB2	1:E:173:SER:HA	2.02	0.40
1:A:97:GLY:HA3	1:A:100:PHE:CD2	2.57	0.40
1:A:35:HIS:CG	1:A:100(C):MET:HE2	2.56	0.40
1:E:11:LEU:HB2	1:E:147:PRO:HG3	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	220/228 (96%)	216 (98%)	3 (1%)	1 (0%)	24	19
1	C	221/228 (97%)	217 (98%)	3 (1%)	1 (0%)	24	19
1	E	220/228 (96%)	216 (98%)	3 (1%)	1 (0%)	24	19
1	H	220/228 (96%)	215 (98%)	5 (2%)	0	100	100
2	B	211/214 (99%)	204 (97%)	7 (3%)	0	100	100
2	D	211/214 (99%)	202 (96%)	9 (4%)	0	100	100
2	F	212/214 (99%)	204 (96%)	8 (4%)	0	100	100
2	L	212/214 (99%)	204 (96%)	8 (4%)	0	100	100
All	All	1727/1768 (98%)	1678 (97%)	46 (3%)	3 (0%)	43	42

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	96	GLY
1	E	96	GLY
1	A	99	GLY

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	184/190 (97%)	177 (96%)	7 (4%)	29	28
1	C	185/190 (97%)	176 (95%)	9 (5%)	22	18

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	184/190 (97%)	177 (96%)	7 (4%)	29	28
1	H	184/190 (97%)	178 (97%)	6 (3%)	33	33
2	B	188/189 (100%)	182 (97%)	6 (3%)	34	35
2	D	188/189 (100%)	184 (98%)	4 (2%)	47	51
2	F	189/189 (100%)	186 (98%)	3 (2%)	55	61
2	L	189/189 (100%)	184 (97%)	5 (3%)	40	42
All	All	1491/1516 (98%)	1444 (97%)	47 (3%)	34	35

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	ARG
1	A	64	LYS
1	A	95	TRP
1	A	150	VAL
1	A	169	VAL
1	A	193	THR
1	A	199	ASN
2	B	19	VAL
2	B	20	THR
2	B	23	CYS
2	B	65	SER
2	B	66	ARG
2	B	129	THR
1	C	64	LYS
1	C	93	SER
1	C	94	ARG
1	C	98	ASP
1	C	148	GLU
1	C	150	VAL
1	C	169	VAL
1	C	193	THR
1	C	199	ASN
2	D	2	ILE
2	D	19	VAL
2	D	20	THR
2	D	129	THR
1	E	64	LYS
1	E	93	SER
1	E	94	ARG

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Mol	Chain	Res	Type
1	E	130	SER
1	E	169	VAL
1	E	193	THR
1	E	199	ASN
2	F	19	VAL
2	F	20	THR
2	F	105	GLU
1	H	64	LYS
1	H	94	ARG
1	H	98	ASP
1	H	169	VAL
1	H	193	THR
1	H	199	ASN
2	L	2	ILE
2	L	19	VAL
2	L	20	THR
2	L	100	GLN
2	L	105	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	ASN
2	B	79	GLN
2	B	91	HIS
1	C	39	GLN
1	C	82(A)	ASN
2	D	38	GLN
2	D	79	GLN
2	D	189	HIS
2	D	210	ASN
2	F	3	GLN
2	F	30	ASN
2	F	91	HIS
1	H	39	GLN
1	H	82(A)	ASN
1	H	164	HIS
2	L	38	GLN
2	L	79	GLN
2	L	152	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	220/228 (96%)	-1.43	0 100 100	17, 30, 61, 75	2 (0%)
1	C	221/228 (96%)	-1.49	0 100 100	15, 29, 56, 74	2 (0%)
1	E	220/228 (96%)	-1.41	0 100 100	14, 35, 59, 75	2 (0%)
1	H	220/228 (96%)	-1.37	0 100 100	17, 36, 58, 72	2 (0%)
2	B	212/214 (99%)	-1.44	0 100 100	18, 37, 48, 60	1 (0%)
2	D	212/214 (99%)	-1.42	0 100 100	21, 37, 54, 62	1 (0%)
2	F	212/214 (99%)	-1.50	0 100 100	17, 34, 49, 57	2 (0%)
2	L	212/214 (99%)	-1.44	0 100 100	17, 34, 59, 70	2 (0%)
All	All	1729/1768 (97%)	-1.44	0 100 100	14, 34, 55, 75	14 (0%)

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.